



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2017 – 08:28 PM EDT

PDB ID : 1JCT
Title : Glucarate Dehydratase, N341L mutant Orthorhombic Form
Authors : Gulick, A.M.; Hubbard, B.K.; Gerlt, J.A.; Rayment, I.
Deposited on : 2001-06-11
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029077
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029077

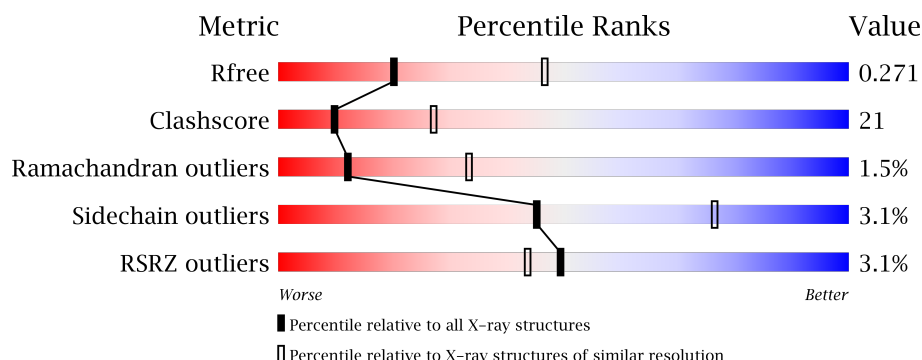
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	3666 (2.80-2.72)
Clashscore	112137	4174 (2.80-2.72)
Ramachandran outliers	110173	4103 (2.80-2.72)
Sidechain outliers	110143	4106 (2.80-2.72)
RSRZ outliers	101464	3697 (2.80-2.72)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	446	<div> <div>3%</div> <div>63%</div> <div>34%</div> <div>..</div> </div>
1	B	446	<div> <div>3%</div> <div>62%</div> <div>35%</div> <div>..</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6935 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Glucarate Dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	0	0
			3421	2163	600	637	21			
1	B	442	Total	C	N	O	S	0	0	0
			3380	2135	592	632	21			

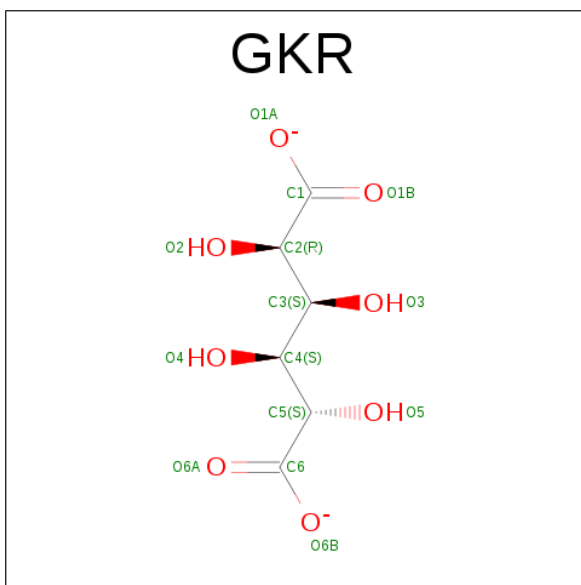
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	341	LEU	ASN	ENGINEERED	UNP P76637
B	341	LEU	ASN	ENGINEERED	UNP P76637

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

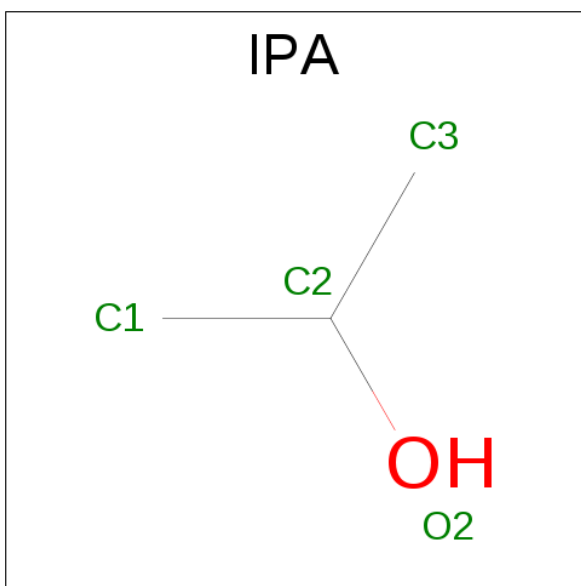
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is D-GLUCARATE (three-letter code: GKR) (formula: C₆H₈O₈).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			14	6	8		
3	B	1	Total	C	O	0	0
			14	6	8		

- Molecule 4 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C_3H_8O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	3	1		
4	B	1	Total	C	O	0	0
			4	3	1		

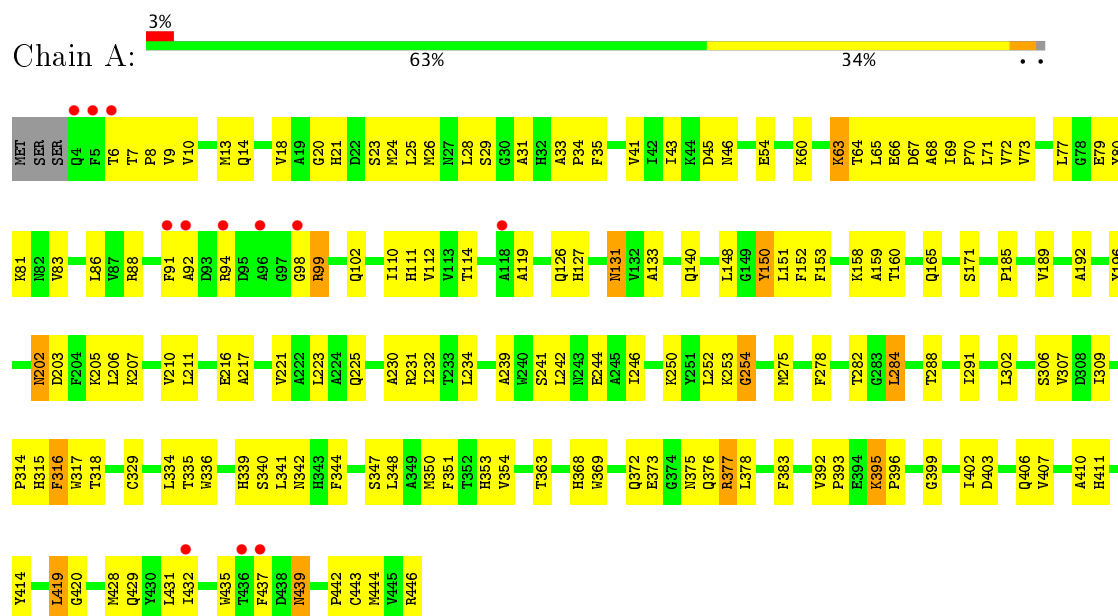
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	62	Total 62	O 62	0	0
5	B	34	Total 34	O 34	0	0

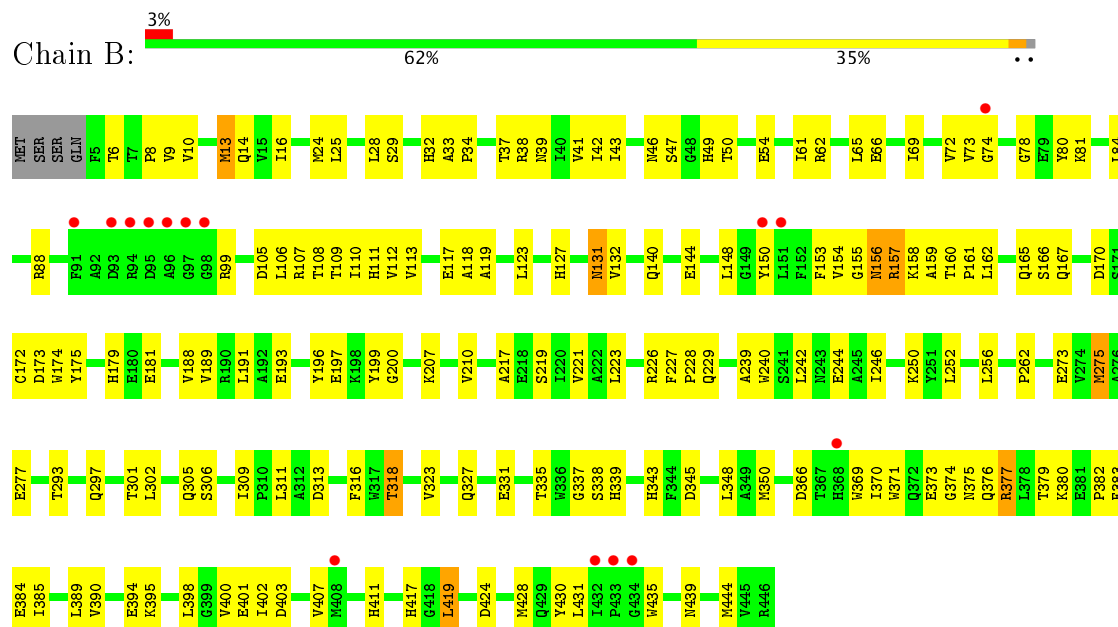
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Glucarate Dehydratase



• Molecule 1: Glucarate Dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	73.92Å 203.16Å 136.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.75 29.19 – 2.75	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.75) 91.0 (29.19-2.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.04 (at 2.76Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.213 , 0.274 0.211 , 0.271	Depositor DCC
R_{free} test set	1182 reflections (4.80%)	DCC
Wilson B-factor (Å ²)	53.8	Xtriage
Anisotropy	0.665	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6935	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, GKR, IPA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.38	0/3500	0.63	0/4743
1	B	0.37	0/3459	0.60	0/4696
All	All	0.38	0/6959	0.61	0/9439

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3421	0	3348	136	0
1	B	3380	0	3258	151	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	14	0	8	0	0
3	B	14	0	8	0	0
4	A	4	0	8	1	0
4	B	4	0	8	3	0
5	A	62	0	0	4	0
5	B	34	0	0	1	0
All	All	6935	0	6638	287	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 21.

All (287) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:ALA:HB1	1:B:252:LEU:HD21	1.51	0.93
1:A:395:LYS:HD2	1:A:399:GLY:HA2	1.55	0.86
1:B:172:CYS:SG	1:B:175:TYR:HD2	1.99	0.86
1:B:43:ILE:HG22	1:B:123:LEU:HD11	1.61	0.82
1:A:275:MET:HG3	1:A:291:ILE:HD13	1.61	0.80
1:A:10:VAL:HG12	1:A:73:VAL:HA	1.63	0.80
1:B:417:HIS:HB2	1:B:419:LEU:HD22	1.66	0.77
1:B:156:ASN:HB2	1:B:181:GLU:OE1	1.83	0.77
1:B:331:GLU:HG2	5:B:610:HOH:O	1.82	0.77
1:B:105:ASP:OD1	1:B:107:ARG:HG3	1.85	0.76
1:A:428:MET:HE1	1:A:443:CYS:HB2	1.67	0.76
1:A:302:LEU:HD22	4:A:601:IPA:H2	1.69	0.74
1:B:13:MET:HG3	1:B:43:ILE:HG13	1.68	0.74
1:B:41:VAL:HG22	1:B:65:LEU:HD21	1.70	0.73
1:A:221:VAL:O	1:A:225:GLN:HG3	1.88	0.73
1:B:16:ILE:HG23	1:B:411:HIS:HD2	1.54	0.72
1:B:293:THR:H	1:B:297:GLN:NE2	1.88	0.71
1:B:61:ILE:HG12	1:B:108:THR:HG23	1.73	0.71
1:B:38:ARG:HD2	1:B:343:HIS:HA	1.73	0.70
1:B:156:ASN:ND2	1:B:159:ALA:HB2	2.07	0.70
1:B:228:PRO:HG2	1:B:229:GLN:HE21	1.56	0.69
1:B:43:ILE:HD13	1:B:119:ALA:CB	2.23	0.69
1:B:189:VAL:O	1:B:193:GLU:HG3	1.93	0.68
1:A:207:LYS:HA	1:A:207:LYS:HE3	1.76	0.68
1:A:316:PHE:O	1:A:316:PHE:HD2	1.78	0.67
1:A:98:GLY:O	1:A:99:ARG:HB2	1.94	0.67
1:A:69:ILE:CG2	1:A:70:PRO:HD3	2.24	0.67
1:A:217:ALA:O	1:A:221:VAL:HG23	1.94	0.67
1:A:71:LEU:HD22	1:A:86:LEU:HD21	1.77	0.67
1:B:10:VAL:CG1	1:B:73:VAL:HA	2.25	0.67
1:A:437:PHE:HE1	1:A:439:ASN:HB3	1.61	0.66
1:A:185:PRO:O	1:A:189:VAL:HG23	1.97	0.65
1:A:6:THR:HG22	1:A:7:THR:N	2.12	0.65
1:B:10:VAL:HG11	1:B:73:VAL:HA	1.80	0.64
1:A:133:ALA:H	1:A:353:HIS:HD2	1.45	0.64
1:B:380:LYS:O	1:B:382:PRO:HD3	1.97	0.64
1:A:432:ILE:O	1:A:435:TRP:HB2	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:GLN:HE22	1:B:375:ASN:HD21	1.47	0.63
1:B:217:ALA:O	1:B:221:VAL:HG23	1.99	0.63
1:A:148:LEU:O	1:A:148:LEU:HD12	1.99	0.62
1:A:60:LYS:HB3	5:A:660:HOH:O	2.00	0.62
1:B:10:VAL:CG1	1:B:74:GLY:H	2.13	0.62
1:A:402:ILE:HD11	1:A:407:VAL:CG2	2.30	0.61
1:B:210:VAL:O	1:B:444:MET:HB2	2.00	0.61
1:A:60:LYS:HE2	1:A:94:ARG:HB3	1.82	0.61
1:A:10:VAL:CG1	1:A:73:VAL:HA	2.30	0.61
1:A:69:ILE:HG22	1:A:70:PRO:HD3	1.83	0.61
1:B:196:TYR:O	1:B:200:GLY:HA2	2.01	0.61
1:A:196:TYR:CD2	1:A:196:TYR:C	2.73	0.61
1:B:193:GLU:O	1:B:196:TYR:HB3	2.01	0.60
1:B:252:LEU:HD12	1:B:256:LEU:HD11	1.83	0.60
1:A:66:GLU:O	1:A:69:ILE:HG22	2.01	0.60
1:A:6:THR:HG22	1:A:7:THR:H	1.67	0.60
1:B:156:ASN:HD22	1:B:159:ALA:HB2	1.67	0.60
1:B:10:VAL:HG13	1:B:74:GLY:H	1.67	0.59
1:A:395:LYS:CD	1:A:399:GLY:HA2	2.27	0.59
1:B:65:LEU:CD1	1:B:112:VAL:HG13	2.32	0.59
1:B:240:TRP:HB3	1:B:244:GLU:HG3	1.86	0.58
1:A:34:PRO:HG2	1:A:35:PHE:CD2	2.38	0.58
1:A:428:MET:HE2	1:A:435:TRP:CE2	2.39	0.58
1:B:65:LEU:HD12	1:B:112:VAL:HG13	1.85	0.57
1:B:175:TYR:CE1	1:B:374:GLY:HA3	2.39	0.57
1:A:210:VAL:O	1:A:444:MET:HB2	2.05	0.57
1:B:24:MET:HE2	1:B:155:GLY:HA3	1.85	0.57
1:A:63:LYS:HE3	1:A:63:LYS:HA	1.85	0.57
1:B:172:CYS:CB	1:B:175:TYR:HD2	2.18	0.57
1:B:6:THR:O	1:B:8:PRO:HD3	2.04	0.57
1:B:50:THR:HA	1:B:123:LEU:HD21	1.86	0.57
1:B:380:LYS:C	1:B:382:PRO:HD3	2.24	0.57
1:B:338:SER:O	1:B:366:ASP:HB2	2.05	0.56
1:B:66:GLU:O	1:B:69:ILE:HG12	2.05	0.56
1:A:348:LEU:HA	1:A:351:PHE:CZ	2.40	0.56
1:A:419:LEU:HD23	1:A:419:LEU:H	1.70	0.56
1:B:242:LEU:O	1:B:246:ILE:HG13	2.06	0.56
1:B:175:TYR:HE1	1:B:374:GLY:HA3	1.71	0.56
1:B:29:SER:O	1:B:424:ASP:HB2	2.05	0.56
1:A:402:ILE:HD11	1:A:407:VAL:HG21	1.88	0.55
1:A:41:VAL:HG22	1:A:65:LEU:HD13	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:309:ILE:HG23	1:B:335:THR:O	2.06	0.55
1:A:8:PRO:HD3	1:A:127:HIS:CE1	2.40	0.55
1:A:376:GLN:O	1:A:377:ARG:HB3	2.06	0.55
1:B:41:VAL:CG2	1:B:65:LEU:HD21	2.35	0.55
1:B:39:ASN:HD21	1:B:62:ARG:HB2	1.72	0.55
1:B:109:THR:O	1:B:112:VAL:HB	2.06	0.55
1:B:37:THR:O	1:B:38:ARG:HG2	2.07	0.55
1:A:252:LEU:O	1:A:254:GLY:N	2.36	0.54
1:A:437:PHE:CE1	1:A:439:ASN:HB3	2.41	0.54
1:B:25:LEU:HB2	1:B:32:HIS:CG	2.42	0.54
1:A:29:SER:HA	1:A:437:PHE:CE2	2.42	0.54
1:A:329:CYS:HA	1:A:334:LEU:HB2	1.89	0.54
1:B:16:ILE:CG2	1:B:411:HIS:HD2	2.20	0.54
1:B:385:ILE:HG12	1:B:390:VAL:CG2	2.37	0.54
1:B:43:ILE:CG2	1:B:123:LEU:HD11	2.37	0.53
1:B:157:ARG:HG3	1:B:158:LYS:N	2.23	0.53
1:A:54:GLU:HB3	1:A:347:SER:OG	2.08	0.53
1:A:18:VAL:HB	1:A:410:ALA:HB1	1.90	0.53
1:A:131:ASN:N	1:A:131:ASN:HD22	2.06	0.53
1:A:316:PHE:O	1:A:316:PHE:CD2	2.60	0.53
1:B:377:ARG:HH11	1:B:377:ARG:HG3	1.74	0.53
1:B:14:GLN:OE1	1:B:16:ILE:HD11	2.09	0.53
1:B:273:GLU:O	1:B:277:GLU:HG3	2.08	0.53
1:A:429:GLN:OE1	1:A:429:GLN:HA	2.09	0.53
1:A:372:GLN:O	1:A:376:GLN:HB2	2.09	0.52
1:B:105:ASP:OD1	1:B:106:LEU:N	2.41	0.52
1:B:197:GLU:HA	1:B:197:GLU:OE1	2.09	0.52
1:A:377:ARG:NH1	1:A:406:GLN:HG3	2.25	0.52
1:B:383:PHE:HD2	1:B:383:PHE:N	2.08	0.52
1:A:241:SER:OG	1:A:244:GLU:HG3	2.09	0.52
1:B:10:VAL:HG11	1:B:72:VAL:O	2.10	0.52
1:B:375:ASN:C	1:B:376:GLN:HG3	2.30	0.52
1:A:210:VAL:HG12	1:A:211:LEU:HD23	1.93	0.51
1:A:13:MET:HG3	1:A:43:ILE:HG12	1.92	0.51
1:A:206:LEU:O	1:A:234:LEU:HD12	2.11	0.51
1:B:42:ILE:C	1:B:43:ILE:HD12	2.31	0.51
1:A:24:MET:CE	1:A:431:LEU:HD11	2.40	0.51
1:B:80:TYR:OH	1:B:81:LYS:HE3	2.11	0.51
1:A:442:PRO:O	1:A:446:ARG:NH1	2.44	0.51
1:A:348:LEU:HD12	1:A:351:PHE:CE2	2.46	0.51
1:A:133:ALA:N	1:A:353:HIS:HD2	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:383:PHE:N	1:B:383:PHE:CD2	2.79	0.51
1:A:29:SER:HA	1:A:437:PHE:CD2	2.46	0.50
1:A:8:PRO:HB3	1:A:45:ASP:OD1	2.10	0.50
1:B:293:THR:H	1:B:297:GLN:HE21	1.58	0.50
1:B:38:ARG:HD3	1:B:54:GLU:OE2	2.10	0.50
1:B:262:PRO:HD2	1:B:275:MET:SD	2.52	0.50
1:B:13:MET:HA	1:B:42:ILE:O	2.11	0.50
1:B:153:PHE:HE1	1:B:188:VAL:HG21	1.77	0.50
1:B:345:ASP:HA	1:B:348:LEU:HB3	1.93	0.50
1:B:62:ARG:HH22	1:B:66:GLU:CD	2.13	0.50
1:B:379:THR:HG23	1:B:401:GLU:O	2.11	0.50
1:A:99:ARG:HG2	1:A:99:ARG:HH11	1.77	0.50
1:B:172:CYS:SG	1:B:175:TYR:CD2	2.83	0.50
1:A:26:MET:HE2	1:A:444:MET:HE3	1.94	0.50
1:A:347:SER:O	1:A:350:MET:HB2	2.12	0.50
1:A:314:PRO:HB3	1:A:354:VAL:HG21	1.93	0.50
1:A:393:PRO:HB2	5:A:616:HOH:O	2.12	0.50
1:B:9:VAL:H	1:B:46:ASN:ND2	2.10	0.50
1:A:344:PHE:CD1	1:A:378:LEU:HD11	2.47	0.49
1:A:369:TRP:O	1:A:373:GLU:HB2	2.12	0.49
1:A:8:PRO:HB2	1:A:77:LEU:CD1	2.42	0.49
1:B:8:PRO:HG3	1:B:127:HIS:CD2	2.47	0.49
1:B:99:ARG:CD	1:B:293:THR:HG21	2.42	0.49
1:B:419:LEU:CD2	1:B:419:LEU:H	2.25	0.49
1:A:153:PHE:HE2	1:A:216:GLU:HG2	1.76	0.49
1:A:282:THR:OG1	1:A:284:LEU:HD23	2.12	0.49
1:B:375:ASN:N	1:B:375:ASN:HD22	2.10	0.49
1:A:288:THR:CG2	1:A:307:VAL:HG21	2.43	0.49
1:A:383:PHE:CD1	1:A:383:PHE:N	2.80	0.49
1:B:131:ASN:HD22	1:B:131:ASN:C	2.14	0.49
1:B:377:ARG:HG3	1:B:377:ARG:NH1	2.27	0.49
1:B:403:ASP:O	1:B:407:VAL:HG23	2.13	0.49
1:A:24:MET:HE1	1:A:431:LEU:HD11	1.95	0.48
1:B:160:THR:C	1:B:162:LEU:H	2.16	0.48
1:B:302:LEU:HD22	4:B:602:IPA:H2	1.95	0.48
1:A:383:PHE:HD1	1:A:383:PHE:N	2.11	0.48
1:B:157:ARG:NH1	1:B:166:SER:CB	2.76	0.48
1:B:172:CYS:HB3	1:B:175:TYR:HB2	1.94	0.48
1:B:323:VAL:O	1:B:327:GLN:HG3	2.12	0.48
1:B:379:THR:HG22	1:B:380:LYS:N	2.27	0.48
1:A:250:LYS:HD2	5:A:655:HOH:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:419:LEU:CD2	1:A:419:LEU:H	2.26	0.48
1:A:99:ARG:HG2	1:A:99:ARG:NH1	2.28	0.48
1:B:80:TYR:CZ	1:B:81:LYS:HG3	2.49	0.48
1:A:67:ASP:O	1:A:70:PRO:HD2	2.14	0.48
1:B:154:VAL:O	1:B:181:GLU:HA	2.12	0.48
1:A:65:LEU:HD23	1:A:112:VAL:HG13	1.96	0.47
1:B:157:ARG:HG3	1:B:158:LYS:H	1.78	0.47
1:A:148:LEU:HD12	1:A:148:LEU:C	2.34	0.47
1:A:336:TRP:O	1:A:363:THR:HG22	2.14	0.47
1:A:80:TYR:OH	1:A:81:LYS:HE2	2.14	0.47
1:B:207:LYS:HA	1:B:207:LYS:HE3	1.95	0.47
1:A:288:THR:HG21	1:A:307:VAL:HG21	1.96	0.47
1:A:439:ASN:HD22	1:A:439:ASN:N	2.11	0.47
1:A:71:LEU:HB3	1:A:86:LEU:HD23	1.96	0.47
1:A:79:GLU:O	1:A:83:VAL:HG23	2.15	0.47
1:B:132:VAL:HG21	1:B:398:LEU:HD23	1.96	0.47
1:A:242:LEU:O	1:A:246:ILE:HG13	2.14	0.47
1:A:411:HIS:O	1:A:414:TYR:HB3	2.15	0.47
1:A:309:ILE:HA	1:A:335:THR:O	2.15	0.46
1:A:43:ILE:HD12	1:A:119:ALA:HB3	1.97	0.46
1:B:193:GLU:OE1	1:B:226:ARG:NE	2.45	0.46
1:A:140:GLN:NE2	1:A:392:VAL:HG11	2.31	0.46
1:A:284:LEU:H	1:A:284:LEU:HD23	1.81	0.46
1:A:114:THR:HG21	1:A:315:HIS:HA	1.98	0.46
1:A:31:ALA:N	1:A:428:MET:HG2	2.29	0.46
1:A:31:ALA:H	1:A:428:MET:HG2	1.81	0.46
1:B:43:ILE:HG22	1:B:123:LEU:CD1	2.41	0.46
1:B:219:SER:O	1:B:223:LEU:HB2	2.16	0.45
1:B:157:ARG:NH2	1:B:158:LYS:NZ	2.64	0.45
1:A:431:LEU:C	1:A:432:ILE:HG13	2.37	0.45
1:B:376:GLN:O	1:B:377:ARG:HB3	2.16	0.45
1:B:380:LYS:HB2	1:B:401:GLU:HB2	1.98	0.45
1:A:203:ASP:OD1	1:A:231:ARG:HB2	2.17	0.45
1:A:8:PRO:HB2	1:A:77:LEU:HD11	1.99	0.45
1:B:196:TYR:O	1:B:200:GLY:CA	2.65	0.45
1:B:428:MET:HG3	1:B:435:TRP:CD2	2.51	0.45
1:B:50:THR:CA	1:B:123:LEU:HD21	2.46	0.45
1:A:10:VAL:HG11	1:A:72:VAL:HG12	1.99	0.45
1:B:302:LEU:HB3	4:B:602:IPA:H2	1.98	0.45
1:B:13:MET:SD	1:B:69:ILE:HD13	2.57	0.45
1:A:60:LYS:CE	1:A:94:ARG:HB3	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:LEU:HB2	1:B:32:HIS:ND1	2.32	0.44
1:B:10:VAL:HG22	1:B:73:VAL:HG22	1.99	0.44
1:A:151:LEU:HD12	1:A:223:LEU:HD21	1.99	0.44
1:A:28:LEU:HD22	1:A:239:ALA:HB2	2.00	0.44
1:B:118:ALA:HB2	1:B:350:MET:CE	2.48	0.44
1:B:25:LEU:HD12	1:B:32:HIS:ND1	2.33	0.44
1:A:278:PHE:C	1:A:278:PHE:CD1	2.91	0.44
1:A:317:TRP:O	1:A:318:THR:HG23	2.18	0.44
1:A:33:ALA:HB1	1:A:34:PRO:HD2	1.99	0.44
1:B:199:TYR:HB3	1:B:385:ILE:CG2	2.48	0.44
1:A:419:LEU:CD2	1:A:419:LEU:N	2.81	0.44
1:B:369:TRP:CE3	1:B:369:TRP:HA	2.53	0.44
1:A:282:THR:OG1	1:A:284:LEU:CD2	2.66	0.44
1:A:68:ALA:HB2	1:A:91:PHE:CE2	2.53	0.44
1:B:293:THR:N	1:B:297:GLN:HE21	2.15	0.44
1:A:6:THR:CG2	1:A:7:THR:N	2.79	0.44
1:A:21:HIS:HB2	1:A:165:GLN:NE2	2.33	0.43
1:B:144:GLU:OE2	1:B:389:LEU:HD22	2.17	0.43
1:B:419:LEU:HD23	1:B:419:LEU:H	1.83	0.43
1:B:131:ASN:HB3	1:B:395:LYS:O	2.18	0.43
1:A:202:ASN:O	1:A:230:ALA:CB	2.66	0.43
1:B:10:VAL:HG13	1:B:73:VAL:HA	1.98	0.43
1:A:8:PRO:HB2	1:A:77:LEU:HG	2.01	0.43
1:B:99:ARG:HG3	1:B:105:ASP:CG	2.39	0.43
1:A:153:PHE:CE2	1:A:216:GLU:HG2	2.53	0.43
1:B:161:PRO:HD2	1:B:430:TYR:CD2	2.54	0.43
1:B:16:ILE:HG23	1:B:411:HIS:CD2	2.44	0.43
1:A:151:LEU:HD11	1:A:192:ALA:HB2	2.00	0.43
1:A:21:HIS:CB	1:A:165:GLN:NE2	2.82	0.43
1:A:288:THR:OG1	1:A:291:ILE:HG12	2.18	0.43
1:A:25:LEU:HD22	1:A:152:PHE:HB2	2.00	0.43
1:B:84:LEU:HD11	1:B:117:GLU:HG3	2.00	0.43
1:B:302:LEU:HD22	4:B:602:IPA:C2	2.49	0.43
1:A:13:MET:HG2	1:A:14:GLN:N	2.33	0.43
1:B:380:LYS:HE3	1:B:402:ILE:O	2.18	0.43
1:A:275:MET:HG3	1:A:291:ILE:CD1	2.41	0.42
1:B:110:ILE:HD11	1:B:318:THR:HG23	2.01	0.42
1:A:340:SER:O	1:A:368:HIS:HE1	2.02	0.42
1:B:301:THR:O	1:B:305:GLN:N	2.50	0.42
1:B:174:TRP:CH2	1:B:370:ILE:HD12	2.54	0.42
1:B:311:LEU:HD23	1:B:337:GLY:HA3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:LYS:C	1:A:160:THR:H	2.23	0.42
1:A:23:SER:HA	5:A:634:HOH:O	2.20	0.42
1:A:9:VAL:HB	1:A:46:ASN:OD1	2.19	0.42
1:B:13:MET:HG3	1:B:43:ILE:CG1	2.45	0.42
1:B:384:GLU:O	1:B:390:VAL:HG22	2.20	0.42
1:A:110:ILE:HG23	1:A:111:HIS:N	2.34	0.42
1:B:226:ARG:HG3	1:B:227:PHE:CE1	2.54	0.42
1:A:69:ILE:HG23	1:A:70:PRO:HD3	1.97	0.42
1:B:246:ILE:O	1:B:250:LYS:HG3	2.20	0.42
1:B:199:TYR:HE2	1:B:373:GLU:HG3	1.85	0.42
1:B:293:THR:N	1:B:297:GLN:NE2	2.64	0.42
1:A:68:ALA:HB2	1:A:91:PHE:HE2	1.85	0.41
1:A:69:ILE:N	1:A:70:PRO:CD	2.83	0.41
1:B:252:LEU:HB2	1:B:256:LEU:HD12	2.02	0.41
1:B:33:ALA:HB1	1:B:34:PRO:HD2	2.00	0.41
1:B:339:HIS:HA	1:B:366:ASP:HB2	2.02	0.41
1:B:41:VAL:HG12	1:B:43:ILE:CD1	2.50	0.41
1:B:199:TYR:HD2	1:B:385:ILE:HB	1.84	0.41
1:B:88:ARG:HB2	1:B:113:VAL:HG21	2.02	0.41
1:A:150:TYR:CZ	1:A:205:LYS:HE2	2.55	0.41
1:B:24:MET:HE3	1:B:431:LEU:HD21	2.02	0.41
1:B:379:THR:CG2	1:B:380:LYS:N	2.82	0.41
1:B:99:ARG:HD2	1:B:293:THR:HG21	2.02	0.41
1:A:20:GLY:O	1:A:35:PHE:HA	2.19	0.41
1:A:314:PRO:CB	1:A:354:VAL:HG21	2.51	0.41
1:B:191:LEU:HD21	1:B:371:TRP:HH2	1.85	0.41
1:A:230:ALA:O	1:A:232:ILE:HD12	2.20	0.41
1:A:316:PHE:CE1	1:A:339:HIS:HE1	2.39	0.41
1:B:111:HIS:NE2	1:B:316:PHE:HB3	2.35	0.41
1:B:316:PHE:HD2	1:B:316:PHE:O	2.04	0.41
1:A:88:ARG:O	1:A:92:ALA:HB2	2.20	0.41
1:B:313:ASP:H	1:B:339:HIS:CE1	2.39	0.41
1:A:341:LEU:O	1:A:342:ASN:C	2.58	0.41
1:B:28:LEU:HD22	1:B:239:ALA:HB2	2.03	0.41
1:B:99:ARG:HG2	1:B:99:ARG:NH1	2.36	0.41
1:A:375:ASN:OD1	1:A:376:GLN:HG3	2.20	0.40
1:A:126:GLN:HA	1:A:396:PRO:HB3	2.02	0.40
1:A:403:ASP:OD1	1:A:406:GLN:HG2	2.21	0.40
1:A:64:THR:HG22	1:A:112:VAL:HG11	2.02	0.40
1:B:10:VAL:HG13	1:B:10:VAL:O	2.21	0.40
1:B:140:GLN:HE21	1:B:140:GLN:HB2	1.58	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:LEU:HD12	1:B:32:HIS:CE1	2.56	0.40
1:B:47:SER:O	1:B:49:HIS:ND1	2.53	0.40
1:B:369:TRP:HA	1:B:369:TRP:HE3	1.87	0.40
1:B:165:GLN:HG3	1:B:179:HIS:CE1	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	441/446 (99%)	398 (90%)	36 (8%)	7 (2%)	11	32
1	B	440/446 (99%)	401 (91%)	33 (8%)	6 (1%)	13	35
All	All	881/892 (99%)	799 (91%)	69 (8%)	13 (2%)	12	33

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	ARG
1	A	102	GLN
1	A	253	LYS
1	B	170	ASP
1	A	159	ALA
1	A	254	GLY
1	B	156	ASN
1	B	318	THR
1	B	439	ASN
1	A	377	ARG
1	B	377	ARG
1	A	420	GLY
1	B	78	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	355/362 (98%)	344 (97%)	11 (3%)	45	76
1	B	345/362 (95%)	334 (97%)	11 (3%)	44	75
All	All	700/724 (97%)	678 (97%)	22 (3%)	45	76

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	LYS
1	A	131	ASN
1	A	150	TYR
1	A	171	SER
1	A	202	ASN
1	A	284	LEU
1	A	306	SER
1	A	316	PHE
1	A	395	LYS
1	A	419	LEU
1	A	439	ASN
1	B	13	MET
1	B	131	ASN
1	B	148	LEU
1	B	150	TYR
1	B	157	ARG
1	B	173	ASP
1	B	275	MET
1	B	306	SER
1	B	394	GLU
1	B	400	VAL
1	B	419	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	39	ASN
1	A	89	ASN
1	A	131	ASN
1	A	141	GLN
1	A	339	HIS
1	A	353	HIS
1	A	372	GLN
1	A	439	ASN
1	B	39	ASN
1	B	82	ASN
1	B	102	GLN
1	B	131	ASN
1	B	140	GLN
1	B	165	GLN
1	B	167	GLN
1	B	179	HIS
1	B	229	GLN
1	B	243	ASN
1	B	297	GLN
1	B	372	GLN
1	B	411	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GKR	A	499	2	7,13,13	2.01	1 (14%)	10,18,18	1.37	1 (10%)
4	IPA	A	601	-	3,3,3	0.55	0	3,3,3	0.12	0
3	GKR	B	502	2	7,13,13	1.88	1 (14%)	10,18,18	1.38	1 (10%)
4	IPA	B	602	-	3,3,3	0.63	0	3,3,3	0.18	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GKR	A	499	2	-	0/12/20/20	0/0/0/0
4	IPA	A	601	-	-	0/0/0/0	0/0/0/0
3	GKR	B	502	2	-	0/12/20/20	0/0/0/0
4	IPA	B	602	-	-	0/0/0/0	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	499	GKR	O2-C2	4.83	1.52	1.42
3	B	502	GKR	O2-C2	4.93	1.52	1.42

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	499	GKR	O5-C5-C4	3.31	114.86	107.52
3	B	502	GKR	O5-C5-C4	3.52	115.32	107.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601	IPA	1	0
4	B	602	IPA	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	443/446 (99%)	-0.07	12 (2%) 55 49	25, 46, 71, 93	0
1	B	442/446 (99%)	0.08	15 (3%) 46 40	28, 60, 84, 97	0
All	All	885/892 (99%)	0.00	27 (3%) 49 44	25, 53, 81, 97	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	96	ALA	5.0
1	B	98	GLY	4.2
1	A	96	ALA	3.7
1	A	94	ARG	2.8
1	B	433	PRO	2.7
1	B	408	MET	2.6
1	A	437	PHE	2.6
1	B	94	ARG	2.5
1	A	118	ALA	2.4
1	B	93	ASP	2.4
1	A	98	GLY	2.3
1	B	432	ILE	2.3
1	A	6	THR	2.2
1	A	4	GLN	2.2
1	A	432	ILE	2.2
1	B	368	HIS	2.1
1	B	97	GLY	2.1
1	B	74	GLY	2.1
1	A	436	THR	2.1
1	B	150	TYR	2.1
1	A	92	ALA	2.1
1	B	95	ASP	2.1
1	A	5	PHE	2.1
1	B	434	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	91	PHE	2.1
1	B	151	LEU	2.0
1	A	91	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	GKR	B	502	14/14	0.94	0.35	1.72	49,57,59,61	0
4	IPA	A	601	4/4	0.93	0.20	1.10	55,58,58,58	0
2	MG	A	498	1/1	0.96	0.21	0.14	35,35,35,35	0
3	GKR	A	499	14/14	0.95	0.21	0.07	40,43,46,47	0
2	MG	B	501	1/1	0.99	0.20	-0.45	29,29,29,29	0
4	IPA	B	602	4/4	0.90	0.29	-	54,55,55,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.